## **Amendments to the Claims**

This listing of claims will replace all prior versions, and listings, of claims in the application:

1. (currently amended) A compound of the formula according to formula (I) below, having the structure:

Formula (I)

wherein:

 $R^1$  is selected from the group consisting of hydrogen, halogen,  $C_{1-4}$  alkyl,  $C_{1-4}$  alkanoyl and aroyl;

G<sup>1</sup> represents is CH<sub>2</sub>-CH<sub>2</sub> or CH=CH;

G<sup>2</sup> represents C<sub>4-7</sub>alkyl or a group of the formula (a), (b) or (c):

X represents a group of the formula (d), (e) or (f):

wherein

Y is, independently, selected from the group consisting of a bond, O, S,  $NR^2$ , -  $NR^2$ -  $C_{1-4}$  alkyl- , and  $C_{1-4}$  alkyl- each of these alkyl groups may contain a heteroatom selected from O,  $NR^2$ , or S;

R<sup>2</sup> represents a is hydrogen, C<sub>1-4</sub> alkyl or C<sub>1-4</sub> alkyl-aryl;

Ar , Ar<sup>4</sup> and Ar<sup>2</sup> are, independently, is selected from the group consisting of an optionally substituted phenyl rings, optionally substituted 5- or 6- membered aromatic heterocyclic rings or <u>an</u> optionally substituted, fused bicyclic, or tricyclic aromatic or heteroaromatic ring systems;

L is, independently, selected from the group consisting of a bond, optionally substituted  $C_{1-4}$  alkanoyl,  $C_{1-4}$  alkenyl,  $C_{1-4}$  alkynyl or  $C_{1-18}$  alkyl which may contain between 0 and 3 heteroatoms independently selected from O,  $NR^2$ , or S; Also, atoms within L may be joined to form up to 3 rings and additionally L may be have up to 3 aryl, heteroaryl or  $-CO_2R^2$  substituents; Z represents a bond, O,  $NR^2$ , S,  $C_{1-4}$  alkylidene or  $C_{1-4}$  alkyl.

- 2. (currently amended) A compound according to claim 1 selected from the group consisting of:
- N-{4-[2-(-[(1R,8S)-11-azatricyclo[6.2.1.02,7]undeca-2,4,6-trien-11-yl)-ethyl]-cyclohexyl}-2-[2-({4-[2-(-[(1R,8S)-11-azatricyclo[6.2.1.02,7]undeca-2,4,6-trien-11-yl)-ethyl]-cyclohexylcarbamoyl}-methyl)-phenyl]-acetamide;
- 2,2'-benzene-1,3-diylbis[N-(trans-4-{2-[(1R,8S)-11-azatricyclo[6.2.1.0 $^{2,7}$ ]undeca-2,4,6-trien-11-yl]ethyl}cyclohexyl)acetamide];
- N-(trans-4-{2-[(1R,8S)-11-azatricyclo[6.2.1.0<sup>2,7</sup>]undeca-2,4,6-trien-11-yl]ethyl}cyclohexyl)-N'-{4-[({[(trans-4-{2-[(1R,8S)-11-azatricyclo[6.2.1.0<sup>2,7</sup>]undeca-2,4,6-trien-11-yl]ethyl}cyclohexyl)amino]carbonyl}amino)methyl]phenyl}urea;
- N',N'''-(benzene-1,3-diyldimethanediyl)bis[N-(trans-4-{2-[(1R,8S)-11-azatricyclo[6.2.1.0<sup>2,7</sup>]undeca-2,4,6-trien-11-yl]ethyl}cyclohexyl)urea];
- N',N'''-1,2-propanediylbis[N-(trans-4-{2-[(1R,8S)-11-azatricyclo[6.2.1.0<sup>2,7</sup>]undeca-2,4,6-trien-11-yl]ethyl}cyclohexyl)urea];
- N',N'''-1,6-hexanediylbis[N-(trans-4-{2-[(1R,8S)-11-azatricyclo[6.2.1. $0^{2,7}$ ]undeca-2,4,6-trien-11-yl]ethyl}cyclohexyl)urea];
- N',N'''-[(methylimino)di-2,1-ethanediyl]bis[N-(trans-4-{2-[(1R,8S)-11-azatricyclo[6.2.1.0<sup>2,7</sup>]undeca-2,4,6-trien-11-yl]ethyl}cyclohexyl)urea];
- N',N'''-1,2-butanediylbis[N-(trans-4-{2-[(1R,8S)-11-azatricyclo[6.2.1.0<sup>2,7</sup>]undeca-2,4,6-trien-11-yl]ethyl}cyclohexyl)urea];

- N-(trans-4-{2-[(1R,8S)-11-azatricyclo[6.2.1.0<sup>2,7</sup>]undeca-2,4,6-trien-11-yl]ethyl}cyclohexyl)-N"-{15-[(trans-4-{2-[(1R,8S)-11-azatricyclo[6.2.1.0<sup>2,7</sup>]undeca-2,4,6-trien-11-yl]ethyl}cyclohexyl)amino]-15-oxo-4,7,10-trioxa-14-azapentadec-1-yl}urea;
- N',N'''-1,4-butanediylbis[N-(trans-4-{2-[(1R,8S)-11-azatricyclo[6.2.1.0<sup>2,7</sup>]undeca-2,4,6-trien-11-yl]ethyl}cyclohexyl)urea];
- N',N'''-(thiodi-2,1-ethanediyl)bis[N-(trans-4-{2-[(1R,8S)-11-azatricyclo[6.2.1.0<sup>2,7</sup>]undeca-2,4,6-trien-11-yl]ethyl]cyclohexyl)urea];
- N',N"'-1,2-ethanediylbis[N-(trans-4-{2-[(1R,8S)-11-azatricyclo[6.2.1.0<sup>2,7</sup>]undeca-2,4,6-trien-11-yl]ethyl}cyclohexyl)urea];
- N-(trans-4-{2-[(1R,8S)-11-azatricyclo[6.2.1.0<sup>2,7</sup>]undeca-2,4,6-trien-11-yl]ethyl}cyclohexyl) N'-[3-([(trans-4-{2-[(1R,8S)-11-azatricyclo[6.2.1.0<sup>2,7</sup>]undeca-2,4,6-trien-11-yl]ethyl}cyclohexyl)amino]carbonyl}amino)-2,2-dimethylpropyl]urea;
- N',N'''-1,2-cyclohexanediylbis[N-(trans-4-{2-[(1R,8S)-11-azatricyclo[6.2.1.0<sup>2,7</sup>]undeca-2,4,6-trien-11-yl]ethyl}cyclohexyl)urea];
- 2,2' benzene 1,4 diylbis[N (trans 4 {2 [(1R,8S)-11-azatricyclo[6.2.1.0<sup>2,7</sup>]undeca-2,4,6-trien-11-yl]ethyl}cyclohexyl)acetamide] trifluoroacetate;
- 2,2'-benzene-1,3-diylbis[N-(trans-4-{2-[(1R,8S)-11-azatricyclo[6.2.1.0<sup>2,7</sup>]undeca-2,4,6-trien-11-yl]ethyl}cyclohexyl)acetamide] trifluoroacetate;
- N-(trans-4-{2-[(1R,8S)-11-azatricyclo[6.2.1.0<sup>2,7</sup>]undeca-2,4,6-trien-11-yl]ethyl}cyclohexyl)-N'-{[5-({[2-({[(trans-4-{2-[(1R,8S)-11-azatricyclo[6.2.1.0<sup>2,7</sup>]undeca-2,4,6-trien-11-yl]ethyl}cyclohexyl)amino]carbonyl}amino)ethyl]thio}methyl)-2-furanyl]methyl}urea;
- N',N'''-(2,3-naphthalenediyldimethanediyl)bis[N-(trans-4-{2-[(1R,8S)-11-azatricyclof6.2.1.0<sup>2,7</sup>]undeca-2,4,6-trien-11-yl]ethyl}cyclohexyl)urea];
- N',N'''-1,12-dodecanediylbis[N-(trans-4-{2-[(1R,8S)-11-azatricyclo[6.2.1.0<sup>2,7</sup>]undeca-2,4,6-trien-11-yl]ethyl}cyclohexyl)urea] trifluoroacetate;
- N',N'''-[benzene-1,2-diylbis(oxy-2,1-ethanediyl)]bis[N-(trans-4-{2-[(1R,8S)-11-azatricyclo[6.2.1.0<sup>2,7</sup>]undeca-2,4,6-trien-11-yl]ethyl}cyclohexyl)urea] trifluoroacetate;

- N',N'''-(benzene-1,3-diyldi-2,1-ethanediyl)bis[N-(trans-4-{2-[(1R,8S)-11-azatricyclo[6.2.1.0<sup>2,7</sup>]undeca-2,4,6-trien-11-yl]ethyl}cyclohexyl)urea] trifluoroacetate;
- N',N'''-(benzene-1,4-diyldi-2,1-ethanediyl)bis[N-(trans-4-{2-[(1R,8S)-11-azatricyclo[6.2.1.0<sup>2,7</sup>]undeca-2,4,6-trien-11-yl]ethyl}cyclohexyl)urea} trifluoroacetate:
- N',N'''-(1,3-cyclohexanediyldimethanediyl)bis[N-(trans-4-{2-[(1R,8S)-11-azatricyclo[6.2.1.0<sup>2,7</sup>]undeca-2,4,6-trien-11-yl]ethyl}cyclohexyl)urea} trifluoroacetate:
- N',N'''-1,11-undecanediylbis[N-(trans-4-{2-[(1R,8S)-11-azatricyclo[6.2.1.0<sup>2,7</sup>]undeca-2,4,6-trien-11-yl]ethyl}cyclohexyl)urea] trifluoroacetate;
- N',N''' 1,10-decanediylbis[N-(trans-4-{2-[(1R,8S)-11-azatricyclo[6.2.1.0<sup>2,7</sup>]undeca-2,4,6-trien-11-yl]ethyl}cyclohexyl)urea] trifluoroacetate;
- N',N'''-1,8-octanediylbis[N-(trans-4-{2-[(1R,8S)-11-azatricyclo[6.2.1.0<sup>2,7</sup>]undeca-2,4,6-trien-11-yl]ethyl}cyclohexyl)urea] trifluoroacetate;
- N',N'''-[1,1'-bi(cyclohexyl)-4,4'-diylbis(oxy-2,1-ethanediyl)]bis[N-(trans-4-{2-[(1R,8S)-11-azatricyclo[6.2.1.0<sup>2,7</sup>]undeca-2,4,6-trien-11-yl]ethyl}cyclohexyl)urea];
- N',N''' (dithiodi 3,1-propanediyl)bis[N (trans 4-{2-[(1R,8S)-11-azatricyclo[6.2.1.0<sup>2,7</sup>]undeca-2,4,6-trien-11-yl]ethyl}cyclohexyl)urea] trifluoroacetate and pharmaceutically acceptable salts thereof.
- 3. (Cancelled)
- 4. (Currently amended) A pharmaceutical composition for the treatment of muscarinic acetylcholine receptor mediated diseases comprising a compound according to claim 1 and a pharmaceutically acceptable carrier thereof.
- 5. (Cancelled)
- 6. (Currently amended) A method of treating a muscarinic acetylcholine receptor mediated disease selected from the group consisting of chronic obstructive lung disease, chronic bronchitis, asthma, chronic respiratory obstruction, pulmonary fibrosis, pulmonary emphysema and allergic rhinitis, wherein acetylcholine binds to

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said receptor, comprising administering a safe and an effective amount of a compound according to claim 1.

## 7. (Cancelled)

- 8. (Currently amended) A method according to claim—7 6 wherein administration is via inhalation via the mouth or nose.
- 9. (previously presented) A method according to claim 8 wherein administration is via a medicament dispenser selected from a reservoir dry powder inhaler, a multidose dry powder inhaler or a metered dose inhaler.
- 10. (currently amended) A method according to claim 9 wherein the compound is administered to a human and has a duration of action of 12 hours or more for a 1 mg dose.
- 11. (currently amended) A method according to claim 10 wherein the compound has a duration of action of 24 12 hours or more.

## 12. (cancelled)

- 13. (new) A compound according to Claim 1 wherein G2 is of formula (a).
- 14. (new) A compound according Claim 1 wherein Ar is phenyl.
- 15. (new) A compound according to Claim 14 wherein Y is independently a C<sub>1-4</sub> alkyl.
- 16. (new) A compound according to Claim 15 wherein G<sup>1</sup> is CH=CH.
- 17. (new) A compound according to Claim 1 wherein G<sup>1</sup> is CH=CH.

18. (new) A compound according to Claim 1 wherein Ar is furyl, thienyl, pyrrolyl, oxazolyl, thiazolyl, imidazolyl, oxadiazolyl, thiadiazolyl, pyridyl, triazolyl, triazinyl, pyridazyl, pyrimidinyl, pyrazolyl, isothiazolyl, or isoxazolyl.

19. (new) A compound according to Claim 17 wherein Ar is furyl, thienyl, pyrrolyl, oxazolyl, thiazolyl, imidazolyl, oxadiazolyl, thiadiazolyl, pyridyl, triazolyl, triazinyl, pyridazyl, pyrimidinyl, pyrazolyl, isothiazolyl, or isoxazolyl.

20. (new) A compound according to Claim 1 wherein Ar is naphthyl, indazolyl, indolyl, benzofuranyl, benzothienyl, benzothiazolyl, benzimidazolyl, benzoxazolyl, benzisoxazolyl, benzisothiazolyl, quinolinyl, quinoxolinyl, quinazolinyl, cinnolinyl, isoquinolinyl, pyrazolo[1,5-a]pyrimidyl, pyrrolo[3,2-b]pyridyl, pyrrolo[3,2-c]pyridyl, thieno[3,2-b]thiophenyl, 1,2-dihydro-2-oxo-quinolinyl, 3,4-dihydro-3-oxo-2*H*-benzoxazinyl, or 1,2-dihydro-2-oxo-3*H*-indolyl.

21. (new) A compound according to Claim 17 wherein Ar is naphthyl, indazolyl, indolyl, benzofuranyl, benzothienyl, benzothiazolyl, benzimidazolyl, benzoxazolyl, benzisoxazolyl, benzisothiazolyl, quinolinyl, quinoxolinyl, quinazolinyl, cinnolinyl, isoquinolinyl, pyrazolo[1,5-a]pyrimidyl, pyrrolo[3,2-b]pyridyl, pyrrolo[3,2-c]pyridyl, thieno[3,2-b]thiophenyl, 1,2-dihydro-2-oxo-quinolinyl, 3,4-dihydro-3-oxo-2*H*-benzoxazinyl, or 1,2-dihydro-2-oxo-3*H*-indolyl.

## 22.(new) A compound of the formula:

wherein X is selected from

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23. A pharmaceutical composition comprising a compound according to Claim 22 and a pharmaceutically acceptable carrier.